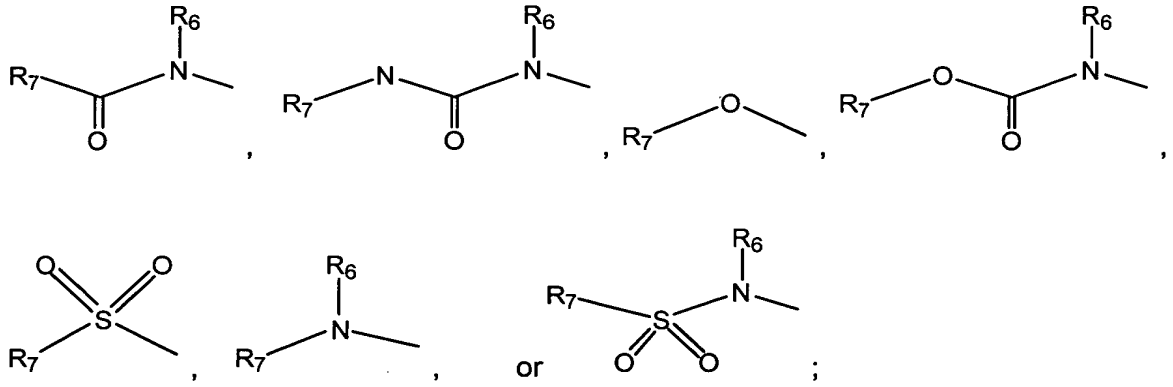
optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_6 alkyl, C_1-C_6 alkoxy, -NO_2, -NH_2, -CN, -CF_3, or -OH;

or a moiety of the formulae:



R_6 is selected from H, C_1-C_6 alkyl, C_1-C_6 alkoxy, -C(O)CH_3, phenyl, -O-phenyl, benzyl, -O-benzyl, the phenyl and benzyl rings of these groups being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_6 alkyl, C_1-C_6 alkoxy, -NO_2, -NH_2, -CN, -CF_3, or -OH;

R_7 is selected from -$(CH_2)_n$-COOH, -$(CH_2)_n$-N-(C_1-C_6 alkyl)$_2$, -$(CH_2)_n$-NH-(C_1-C_6 alkyl), -CF_3, C_1-C_6 alkyl, C_3-C_5 cycloalkyl, C_1-C_6 alkoxy, -NH-(C_1-C_6 alkyl), -N-(C_1-C_6 alkyl)$_2$, pyridinyl, thienyl, furyl, pyrrolyl, quinolyl, $(CH_2)_n$phenyl, phenyl,-O-phenyl, benzyl, -O-benzyl, adamantyl, or morpholinyl, -$(CH_2)_n$-phenyl-O-phenyl, -$(CH_2)_n$-phenyl-CH_2-phenyl, -$(CH_2)_n$-O-phenyl-CH_2-phenyl, -$(CH_2)_n$-phenyl-(O-CH_2-phenyl)$_2$,the rings of these groups being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_6 alkyl, C_1-C_6 alkoxy, -NH_2, -NO_2, -CF_3,CO_2H, or -OH;

R_2 is selected from H, halogen, -CF_3, -OH, -C_1-C_{10} alkyl, C_1-C_{10} alkoxy, -CHO, -CN, -NO_2, -NH_2, -NH-C_1-C_6 alkyl, -N(C_1-C_6 alkyl)$_2$, -N-SO_2-C_1-C_6 alkyl, or -SO_2-C_1-C_6 alkyl;

R_3 is selected from H, -CF_3, C_1-C_6 lower alkyl, C_1-C_6 lower alkoxy, C_3-C_{10} cycloalkyl, -C_1-C_6 alkyl, -C_3-C_{10} cycloalkyl, -CHO, halogen, $(CH_2)_n$C(O)NH_2 or a moiety of the formula – L^1-M^1:

L^1 indicates a linking or bridging group of the formulae -$(CH_2)_n$-, -S-, -O-,

$-C(O)-$, $-(CH_2)_n-C(O)-$, $-(CH_2)_n-C(O)-(CH_2)_n-$, $-(CH_2)_n-O-(CH_2)_n-$, or $-(CH_2)_n-S-(CH_2)_n-$, $C(O)C(O)X$, $-(CH_2)_n-N-(CH_2)_n-$;

M^1 is selected from the group consisting of:

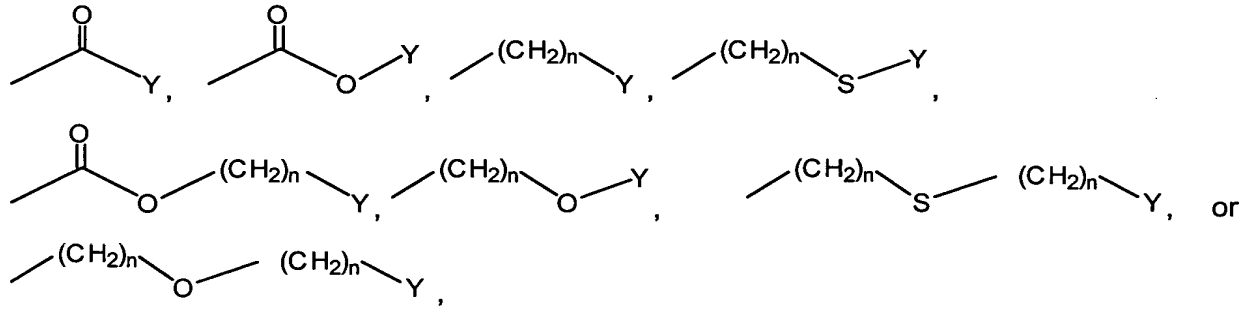
a) H, C_1-C_6 lower alkyl, C_1-C_6 lower alkoxy, C_3-C_{10} cycloalkyl, phenyl, and benzyl, the cycloalkyl, phenyl and benzyl rings being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_{10} alkyl, C_1-C_{10} alkoxy, $-NO_2$, $-NH_2$, -CN, and $-CF_3$, with the proviso that M^1 cannot be H when L^1 is -O-;

b) a six-membered heterocyclic ring containing one, two or three ring heteroatoms selected from N, S or O, the six-membered heterocyclic ring being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_{10} alkyl, C_1-C_{10} alkoxy, -CHO, $-NO_2$, $-NH_2$, -CN, $-CF_3$ or -OH; and

c) a bicyclic ring moiety containing from 8 to 10 ring atoms and optionally containing from 1 to 3 ring heteroatoms selected from N, S or O, the bicyclic ring moiety being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_{10} alkyl, C_1-C_{10} alkoxy, -CHO, $-NO_2$, $-NH_2$, -CN, $-CF_3$ or -OH;

R_4 is selected from the group of C_1-C_6 lower alkyl, C_1-C_6 lower alkoxy, $-(CH_2)_n-C_3-C_6$ cycloalkyl, $-(CH_2)_n-S-(CH_2)_n-C_3-C_5$ cycloalkyl, $-(CH_2)_n-O-(CH_2)_n-C_3-C_5$ cycloalkyl, or the groups of:

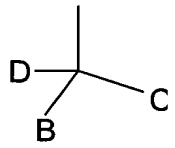
a) $-(CH_2)_n$-phenyl-O-phenyl, $-(CH_2)_n$-phenyl-CH_2-phenyl, $-(CH_2)_n$-O-phenyl-CH_2-phenyl, $-(CH_2)_n$-phenyl-(O-CH_2-phenyl)$_2$, or a moiety of the formulae:



wherein n is independently selected in each appearance as an integer from 0 to 3, Y is C_3-C_5 cycloalkyl, phenyl, benzyl, napthyl, pyridinyl, quinolyl, furyl, thienyl or pyrrolyl; rings of these

groups being optionally substituted by from 1 to 3 substituents selected from H, halogen, -CF_3, -OH, -C_1-C_6 alkyl, C_1-C_6 alkoxy, -NH_2, -NO_2 or a five membered heterocyclic ring containing one heteroatom selected from N, S, or O; or

b) a moiety of the formulae -$(CH_2)_n$-A, -$(CH_2)_n$-S-A, or -$(CH_2)_n$-O-A, wherein A is the moiety:

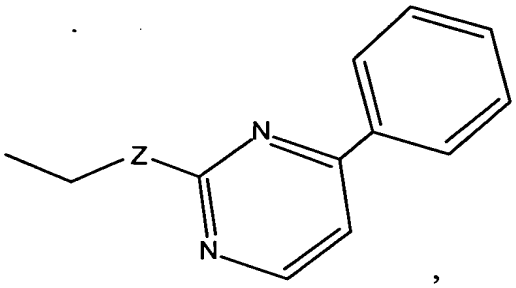


wherein

D is H, C_1-C_6 lower alkyl, C_1-C_6 lower alkoxy, or -CF_3;

B and C are independently selected from phenyl, pyridinyl, furyl, thienyl, pyrimidinyl or pyrrolyl groups, each optionally substituted by from 1 to 3, substituents selected from H, halogen, -CF_3, -OH, -C_1-C_6 alkyl, C_1-C_6 alkoxy, or -NO_2; or

c) a moiety of the formulae:



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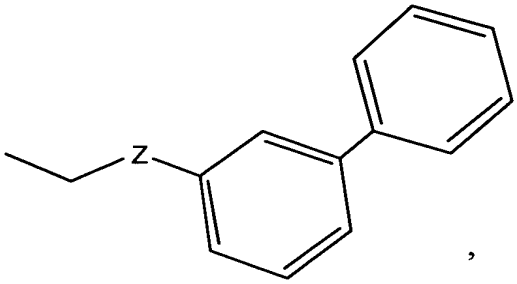

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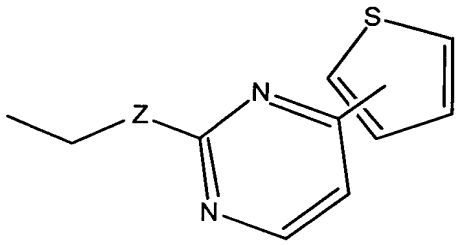

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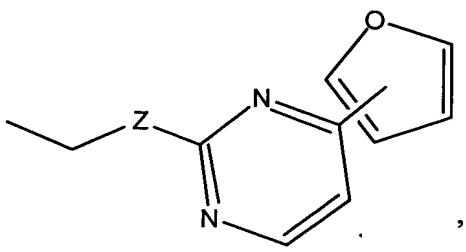

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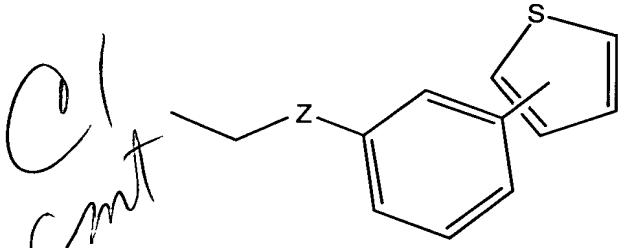
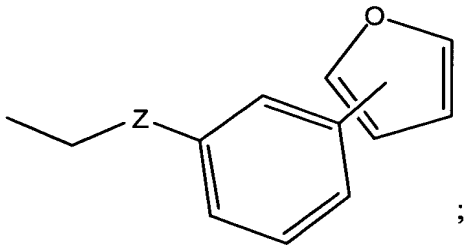


, or


;

wherein Z is O or S and the phenyl and pyrimidinyl rings of each moiety are optionally and independently substituted by from 1 to 3 substituents selected from halogen, $-CF_3$, -OH, $-C_1$-C_6 alkyl, C_1-C_6 alkoxy, $-NH_2$, or $-NO_2$; or

d) a moiety of the formula $-L^2-M^2$, wherein:

L^2 indicates a linking or bridging group of the formulae $-(CH_2)_n-$, -S-, -O-, $-SO_2-$, -C(O)-, $-(CH_2)_n-C(O)-$, $-(CH_2)_n-C(O)-(CH_2)_n-$, $-(CH_2)_n-O-(CH_2)_n-$, or $-(CH_2)_n-S-(CH_2)_n-$, -C(O)C(O)X;
where X = O, N

M^2 is selected from the group of C_1-C_6 lower alkyl, C_1-C_6 lower alkoxy, C_3-C_{10} cycloalkyl, phenyl or benzyl, the cycloalkyl, phenyl or benzyl rings being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_{10} alkyl, C_1-C_{10} alkoxy, $-NO_2$, $-NH_2$, -CN, or $-CF_3$; or

i) a five-membered heterocyclic ring containing one or two ring heteroatoms selected from N, S or O, the five-membered heterocyclic ring being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_{10} alkyl, C_1-C_{10} alkoxy, $-NO_2$, $-NH_2$, -CN, or $-CF_3$; or

ii) a six-membered heterocyclic ring containing one, two or three ring heteroatoms selected from N, S or O, the six-membered heterocyclic ring being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_{10} alkyl, C_1-C_{10} alkoxy, -CHO, $-NO_2$, $-NH_2$, -CN, $-CF_3$ or -OH; or

iii) a bicyclic ring moiety containing from 8 to 10 ring atoms and optionally containing from 1 to 3 ring heteroatoms selected from N, S or O, the bicyclic ring moiety being optionally substituted by from 1 to 3 substituents selected from halogen, C_1-C_{10} alkyl, C_1-C_{10} alkoxy, -CHO, $-NO_2$, $-NH_2$, -CN, $-CF_3$ or -OH;

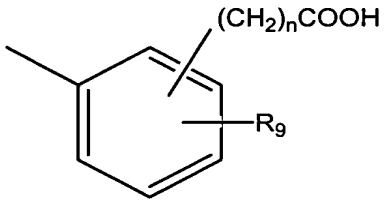
Cont

n is an integer from 0 to 3;

R_5 is a moiety selected from the formulae $-L^3$-M^3

wherein L^3 is a bridging or linking moiety selected from a chemical bond, $-(CH_2)_n-$, -S-, -O-, $-SO_2-$, -C(O)-, $-(CH_2)_n$-C(O)-, $-(CH_2)_n$-C(O)-$(CH_2)_n-$, $-(CH_2)_n$-O-$(CH_2)_n-$,$-(CH_2)_n$-S-$(CH_2)_n-$, -C(Z)-N(R_6)-, -C(Z)-N(R_6)-$(CH_2)_n-$, -C(O)-C(Z)-N(R_6)-, -C(O)-C(Z)-N(R_6)-$(CH_2)_n-$, -C(Z)-NH-SO_2-, -C(Z)-NH-SO_2-$(CH_2)_n-$, $-(CH_2)_n$-S-$(CH_2)_n-$, $-(CH_2)_n$-SO-$(CH_2)_n-$, $-(CH_2)_n-SO_2-(CH_2)_n-$, or $-(CH_2)_n$-CH=CH-$(CH_2)_n$-O-;

M^3 is



and n is an integer from 0 to 3;

R_9 is selected from H, halogen, $-CF_3$, -OH, -COOH, $-(CH_2)_n$-COOH, $-(CH_2)_n$-C(O)-COOH, $-C_1$-C_6 alkyl, -O-C_1-C_6 alkyl, -NH(C_1-C_6 alkyl), or -N(C_1-C_6 alkyl)$_2$;

n is an integer from 0 to 3;

or a pharmaceutically acceptable salt thereof.